NEWS LOGIN

NEWS IPC8

Welcome to STN International! Enter x:x LOGINID: SSPTADEG1625 PASSWORD: TERMINAL (ENTER 1, 2, 3, OR ?):2 \* \* \* \* \* \* \* \* \* \* Welcome to STN International Web Page for STN Seminar Schedule - N. America NEWS NEWS MAR 31 IFICDB, IFIPAT, and IFIUDB enhanced with new custom IPC display formats MAR 31 NEWS 3 CAS REGISTRY enhanced with additional experimental spectra NEWS MAR 31 CA/CAplus and CASREACT patent number format for U.S. applications updated NEWS MAR 31 LPCI now available as a replacement to LDPCI NEWS MAR 31 EMBASE, EMBAL, and LEMBASE reloaded with enhancements 6 7 NEWS APR 04 STN AnaVist, Version 1, to be discontinued WPIDS, WPINDEX, and WPIX enhanced with new NEWS 8 APR 15 predefined hit display formats NEWS 9 APR 28 EMBASE Controlled Term thesaurus enhanced NEWS 10 APR 28 IMSRESEARCH reloaded with enhancements NEWS 11 MAY 30 INPAFAMDB now available on STN for patent family searching DGENE, PCTGEN, and USGENE enhanced with new homology NEWS 12 MAY 30 sequence search option NEWS 13 JUN 06 EPFULL enhanced with 260,000 English abstracts NEWS 14 JUN 06 KOREAPAT updated with 41,000 documents NEWS 15 JUN 13 USPATFULL and USPAT2 updated with 11-character patent numbers for U.S. applications NEWS 16 JUN 19 CAS REGISTRY includes selected substances from web-based collections NEWS 17 JUN 25 CA/CAplus and USPAT databases updated with IPC reclassification data NEWS 18 JUN 30 AEROSPACE enhanced with more than 1 million U.S. patent records NEWS 19 JUN 30 EMBASE, EMBAL, and LEMBASE updated with additional options to display authors and affiliated organizations NEWS 20 JUN 30 STN on the Web enhanced with new STN AnaVist Assistant and BLAST plug-in JUN 30 NEWS 21 STN AnaVist enhanced with database content from EPFULL NEWS 22 JUL 28 CA/CAplus patent coverage enhanced NEWS 23 JUL 28 EPFULL enhanced with additional legal status information from the epoline Register JUL 28 IFICDB, IFIPAT, and IFIUDB reloaded with enhancements NEWS 24 NEWS 25 JUL 28 STN Viewer performance improved NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3, AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008. NEWS HOURS STN Operating Hours Plus Help Desk Availability

Welcome Banner and News Items

For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 17:45:17 ON 28 JUL 2008

=> file registry COST IN U.S. DOLLARS

SINCE FILE TOTAL
ENTRY SESSION
0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 17:45:40 ON 28 JUL 2008
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
COPYRIGHT (C) 2008 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 27 JUL 2008 HIGHEST RN 1036536-16-9 DICTIONARY FILE UPDATES: 27 JUL 2008 HIGHEST RN 1036536-16-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH January 9, 2008.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

=>

Uploading C:\Program Files\Stnexp\Queries\10-518405genBb.str

chain nodes :

14 15 16 17 18 19 20 22 23

ring nodes :

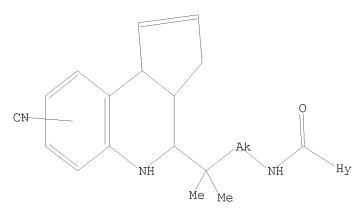
1 2 3 4 5 6 7 8 9 10 11 12 13 chain bonds: 9-14 14-15 14-22 14-23 15-16 16-17 17-18 17-19 ring bonds: 1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 7-11 8-9 8-13 9-10 11-12 12-13 exact/norm bonds: 5-7 6-10 7-8 7-11 8-9 8-13 9-10 11-12 12-13 14-15 15-16 16-17 17-18 17-19 exact bonds: 9-14 14-22 14-23 normalized bonds: 1-2 1-6 2-3 3-4 4-5 5-6

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 15:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:Atom 22:CLASS 23:CLASS

## L1 STRUCTURE UPLOADED

=> d 11 L1 HAS NO ANSWERS L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss full FULL SEARCH INITIATED 17:46:15 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 621 TO ITERATE

100.0% PROCESSED 621 ITERATIONS 13 ANSWERS SEARCH TIME: 00.00.01

L2 13 SEA SSS FUL L1

=> file caplus
COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST
178.36
178.57

FILE 'CAPLUS' ENTERED AT 17:46:25 ON 28 JUL 2008

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2008 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December 26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 28 Jul 2008 VOL 149 ISS 5 FILE LAST UPDATED: 27 Jul 2008 (20080727/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

http://www.cas.org/legal/infopolicy.html

=> s 12

L3 1 L2

=> d 13 abs ibib hitstr

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN GI

AB The title nonsteroidal tetrahydroquinoline derivs. with general formula of I [wherein R1 = NO2 or CN; X = CH or O; m = 0 or 1; Y = (un)substituted alkylene; R2 = H, alkyl, cycloalkyl, or aralkyl; Z = (un)substituted alkyl, aryl, etc.] or pharmaceutically acceptable salts thereof are prepared as androgen receptor agonists. For example, the compound II was prepared in a three-step synthesis starting from 4-nitroaniline, cyclopentadiene, and tert-Bu N-(2,2,-dimethyl-3-oxopropyl)carbamate. II showed relative binding affinity of 1076 against androgen receptor in rat. Formulations containing I as an active ingredient were also described.

ACCESSION NUMBER: 2004:2862 CAPLUS

DOCUMENT NUMBER: 140:59527

TITLE: Preparation of bicyclic tetrahydroquinoline

derivatives as androgen receptor agonists

INVENTOR(S): Miyakawa, Motonori; Sumita, Yuji; Furuya, Kazuyuki;

Ichikawa, Kiyonoshin; Yamamoto, Noriko; Hanada, Keigo;

Amano, Seiji; Nejishima, Hiroaki

PATENT ASSIGNEE(S): Kaken Pharmaceutical Co., Ltd., Japan

PCT Int. Appl., 85 pp. SOURCE:

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

	PAT						KIND DATE			APPLICATION NO.								
	WO	0 2004000816								WO 2003-JP7799								
		W:	ΑE,	AG,	AL,	AM,	ΑT,	ΑU,	ΑZ,	BA,	BB,	BG,	BR,	BY,	BZ,	CA,	CH,	CN,
			CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	ES,	FI,	GB,	GD,	GE,	GH,
			GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	KE,	KG,	KP,	KR,	KΖ,	LC,	LK,	LR,
			LS,	LT,	LU,	LV,	MA,	MD,	MG,	MK,	MN,	MW,	MX,	MZ,	NI,	NO,	NZ,	OM,
			PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	ΤJ,	TM,	TN,	TR,
			TT,	TZ,	UA,	UG,	US,	UZ,	VC,	VN,	YU,	ZA,	ZM,	ZW				
		RW:	GH,	GM,	ΚE,	LS,	MW,	MZ,	SD,	SL,	SZ,	TZ,	UG,	ZM,	ZW,	ΑM,	ΑZ,	BY,
			KG,	KΖ,	MD,	RU,	ΤJ,	TM,	ΑT,	BE,	BG,	CH,	CY,	CZ,	DE,	DK,	EE,	ES,
			FI,	FR,	GB,	GR,	ΗU,	ΙE,	ΙΤ,	LU,	MC,	NL,	PT,	RO,	SE,	SI,	SK,	TR,
			BF,	ΒJ,	CF,	CG,	CI,	CM,	GΑ,	GN,	GQ,	GW,	ML,	MR,	NE,	SN,	TD,	TG
	AU 2003244313									AU 2003-244313								
	ΕP	1520																
		R:	,		,	,	,	ES,	,	,	,	,	,	,	,	,	,	PT,
								RO,										
		2006						2006	0615		US 2	005-	5184	05		2	0051	118
PRIO	RIT	Z APP	LN.	INFO	.:						JP 2	002-	1790	88		A 2	0020	619
											WO 2	003-	JP'/'/	99	1	W 2	0030	619
		DURCE																
ΙT		7333-								-								
		7333-																
		7333-																
		7333-			/333	-99-	4P 6	3/33	4-00	-0P								
		7334-			7		,			CDM		. 1						
		: PAC				_			_		_							
		nerap	eutı	c us	e);	RIOT	(BI	отод	ıcaı	stu	ay);	PRE	P (P.	repa	ratı	on);	USE	<b>ప</b>
	( U S	ses)		امائتما					-£ 1		_12_		la	al		1 4	al a .a	
			_				-	t10n.	OI ]	отсй	CIIC	tet	rany	ar.od.	ulliO	ттие	aer	ivs.

as androgen receptor agonists)

RN 637333-89-2 CAPLUS

3-Pyridinecarboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-tetCN cyclopenta[c]quinolin-4-y1]-2-methylpropy1]-, rel- (CA INDEX NAME)

RN 637333-90-5 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-methyl-, rel- (CA INDEX NAME)

Relative stereochemistry.

RN 637333-91-6 CAPLUS

CN 3-Pyridinecarboxamide, 6-chloro-N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

Relative stereochemistry.

RN 637333-92-7 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-1,6-dihydro-6-oxo-, rel- (CA INDEX NAME)

RN 637333-93-8 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-methoxy-, rel- (CA INDEX NAME)

Relative stereochemistry.

RN 637333-94-9 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-ethoxy-, rel- (CA INDEX NAME)

RN 637333-95-0 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-(2, 2, 2-trifluoroethoxy)-, rel- (CA INDEX NAME)

Relative stereochemistry.

RN 637333-96-1 CAPLUS

CN 3-Pyridinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-6-(methoxymethoxy)-, rel- (CA INDEX NAME)

RN 637333-97-2 CAPLUS

CN 3-Pyridinecarboxamide, 6-(acetylamino)-N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel-(CA INDEX NAME)

Relative stereochemistry.

RN 637333-98-3 CAPLUS

CN 2-Pyrazinecarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-5-methyl-, rel- (CA INDEX NAME)

RN 637333-99-4 CAPLUS

CN 1H-Pyrrole-2-carboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

Relative stereochemistry.

RN 637334-00-0 CAPLUS

CN 1H-Pyrrole-2-carboxamide, N-[2-[(3aR, 4S, 9bS)-8-cyano-3a, 4, 5, 9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-1-methyl-, rel- (CA INDEX NAME)

RN 637334-01-1 CAPLUS

CN 2-Furancarboxamide, N-[2-[(3aR,4S,9bS)-8-cyano-3a,4,5,9b-tetrahydro-3H-cyclopenta[c]quinolin-4-yl]-2-methylpropyl]-, rel- (CA INDEX NAME)

Relative stereochemistry.

REFERENCE COUNT:

THERE ARE 10 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> log off
ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF
LOGOFF? (Y)/N/HOLD:y
STN INTERNATIONAL LOGOFF AT 17:48:10 ON 28 JUL 2008